



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 14, 2023 – 11:23 PM JST

PDB ID : 6IH8  
Title : Crystal structure of Phosphite Dehydrogenase mutant I151R/P176R/M207A from *Ralstonia* sp. 4506  
Authors : Song, X.; Feng, Y.; Liu, Y.; Zhao, Z.  
Deposited on : 2018-09-28  
Resolution : 2.25 Å (reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

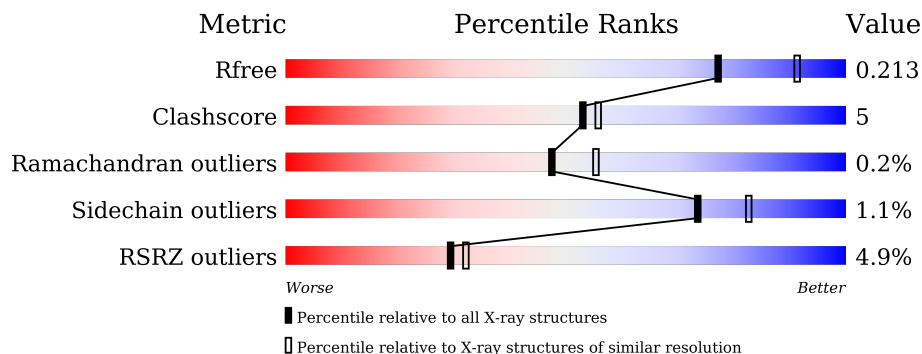
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.25 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	338	<div style="display: flex; align-items: center;"> <div style="width: 4%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 86%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 11%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 0%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="margin-left: 40px;">86% <span style="float: right;">11% •</span></p>
1	B	338	<div style="display: flex; align-items: center;"> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 83%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 14%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 0%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="margin-left: 40px;">83% <span style="float: right;">14% •</span></p>
1	C	338	<div style="display: flex; align-items: center;"> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 85%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 12%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 0%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="margin-left: 40px;">85% <span style="float: right;">12% ••</span></p>
1	D	338	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 85%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 11%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 0%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="margin-left: 40px;">85% <span style="float: right;">11% •</span></p>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10702 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Phosphite dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	329	2525	1604	441	465	15	0	0	0
1	B	329	2525	1604	441	465	15	0	0	0
1	C	330	2533	1609	442	466	16	0	0	0
1	D	326	2501	1589	438	459	15	0	0	0

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	151	ARG	ILE	engineered mutation	UNP G4XDR8
A	176	ARG	PRO	engineered mutation	UNP G4XDR8
A	207	ALA	MET	engineered mutation	UNP G4XDR8
A	337	LEU	-	expression tag	UNP G4XDR8
A	338	GLU	-	expression tag	UNP G4XDR8
B	151	ARG	ILE	engineered mutation	UNP G4XDR8
B	176	ARG	PRO	engineered mutation	UNP G4XDR8
B	207	ALA	MET	engineered mutation	UNP G4XDR8
B	337	LEU	-	expression tag	UNP G4XDR8
B	338	GLU	-	expression tag	UNP G4XDR8
C	151	ARG	ILE	engineered mutation	UNP G4XDR8
C	176	ARG	PRO	engineered mutation	UNP G4XDR8
C	207	ALA	MET	engineered mutation	UNP G4XDR8
C	337	LEU	-	expression tag	UNP G4XDR8
C	338	GLU	-	expression tag	UNP G4XDR8
D	151	ARG	ILE	engineered mutation	UNP G4XDR8
D	176	ARG	PRO	engineered mutation	UNP G4XDR8
D	207	ALA	MET	engineered mutation	UNP G4XDR8
D	337	LEU	-	expression tag	UNP G4XDR8
D	338	GLU	-	expression tag	UNP G4XDR8

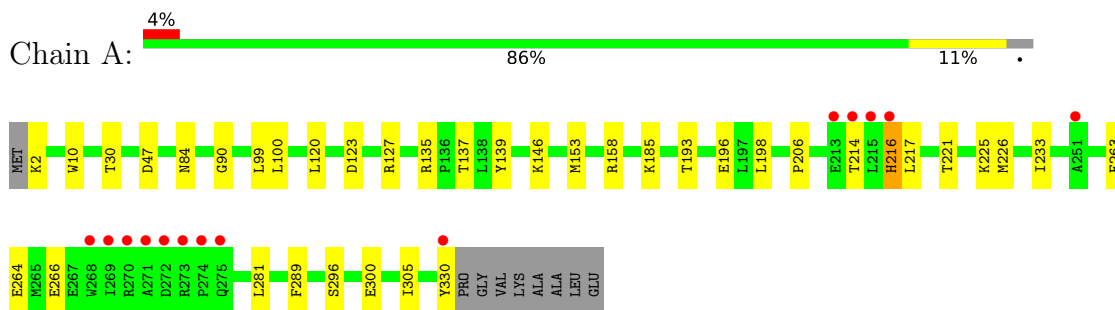
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	161	Total 161	O 161	0	0
2	B	138	Total 138	O 138	0	0
2	C	143	Total 143	O 143	0	0
2	D	176	Total 176	O 176	0	0

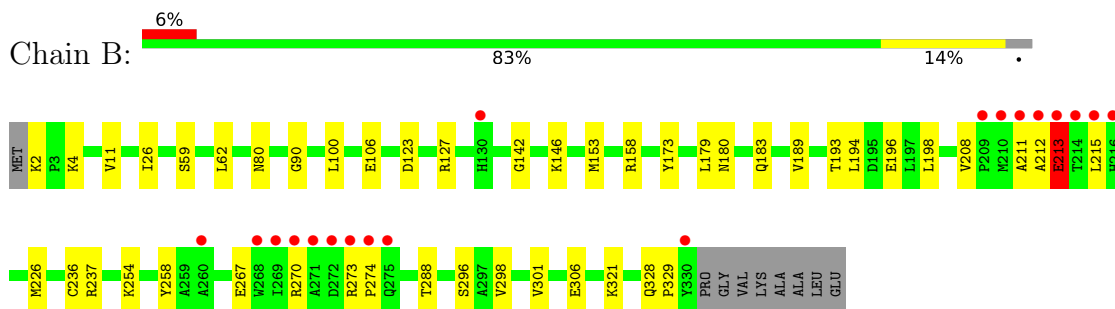
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

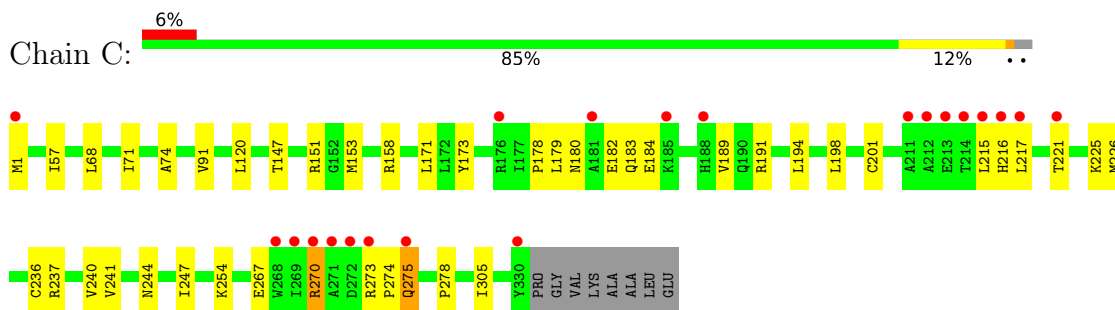
- Molecule 1: Phosphite dehydrogenase



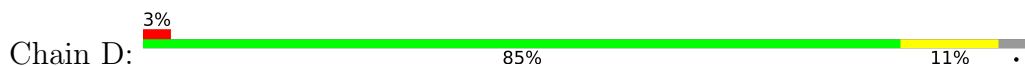
- Molecule 1: Phosphite dehydrogenase

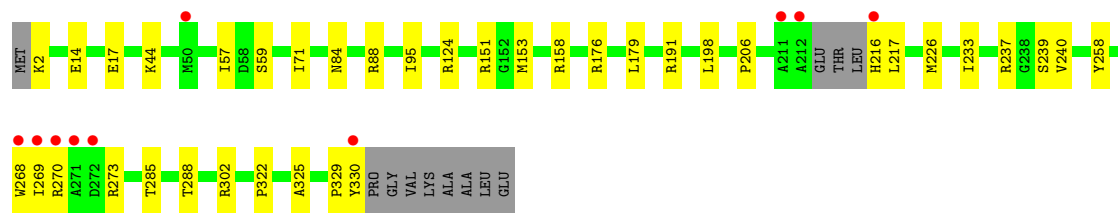


- Molecule 1: Phosphite dehydrogenase



- Molecule 1: Phosphite dehydrogenase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	62.03Å 118.59Å 98.23Å 90.00° 90.65° 90.00°	Depositor
Resolution (Å)	26.75 – 2.25 27.66 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.1 (26.75-2.25) 99.1 (27.66-2.25)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.30 (at 2.24Å)	Xtrriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, $R_{free}$	0.176 , 0.213 0.176 , 0.213	Depositor DCC
$R_{free}$ test set	3261 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.7	Xtrriage
Anisotropy	0.091	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 39.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.097 for h,-k,-l	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10702	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.43% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.43	0/2574	0.57	0/3498
1	B	0.42	0/2574	0.58	0/3498
1	C	0.41	0/2582	0.57	1/3508 (0.0%)
1	D	0.42	0/2549	0.59	0/3462
All	All	0.42	0/10279	0.58	1/13966 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	120	LEU	CA-CB-CG	-5.19	103.36	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	213	GLU	Peptide

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within



the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2525	0	2566	31	0
1	B	2525	0	2566	28	0
1	C	2533	0	2578	29	0
1	D	2501	0	2541	25	0
2	A	161	0	0	6	0
2	B	138	0	0	3	0
2	C	143	0	0	0	0
2	D	176	0	0	3	0
All	All	10702	0	10251	110	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (110) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:214:THR:HB	1:A:217:LEU:HD11	1.42	0.97
1:C:273:ARG:HH22	1:C:275:GLN:HE21	1.17	0.89
1:B:211:ALA:HA	1:D:95:ILE:HD11	1.64	0.79
1:A:217:LEU:N	1:A:217:LEU:HD22	2.02	0.74
1:B:180:ASN:OD1	1:B:183:GLN:N	2.20	0.72
1:B:153:MET:HE2	1:B:179:LEU:HD11	1.70	0.72
1:B:226:MET:O	1:B:254:LYS:NZ	2.22	0.72
1:C:273:ARG:HH12	1:C:275:GLN:HA	1.55	0.72
1:A:153:MET:HE1	1:A:158:ARG:HA	1.72	0.71
1:B:4:LYS:HG3	1:B:26:ILE:HD13	1.77	0.67
1:B:2:LYS:N	2:B:404:HOH:O	2.28	0.66
1:D:216:HIS:N	2:D:404:HOH:O	2.28	0.66
1:C:273:ARG:HH22	1:C:275:GLN:NE2	1.93	0.64
1:B:208:VAL:HG23	1:B:236:CYS:SG	2.38	0.62
1:D:176:ARG:HG2	1:D:176:ARG:HH11	1.62	0.62
1:C:240:VAL:HG23	1:C:241:VAL:HG23	1.82	0.62
1:C:273:ARG:NH1	1:C:274:PRO:O	2.32	0.62
1:A:214:THR:CB	1:A:217:LEU:HD11	2.25	0.61
1:B:153:MET:HE3	1:B:158:ARG:HG2	1.83	0.61
1:A:185:LYS:HG2	2:A:537:HOH:O	2.01	0.60
1:C:153:MET:HE1	1:C:158:ARG:HA	1.83	0.60
1:C:215:LEU:HB2	1:C:216:HIS:HD2	1.66	0.60
1:D:268:TRP:HA	1:D:273:ARG:HH11	1.69	0.58
1:B:11:VAL:HG12	1:B:306:GLU:HG2	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:ASN:ND2	2:A:401:HOH:O	2.17	0.57
1:D:124:ARG:NH2	1:D:285:THR:HG23	2.20	0.56
1:C:273:ARG:NH1	1:C:275:GLN:HA	2.20	0.56
1:C:179:LEU:O	1:C:183:GLN:NE2	2.37	0.56
1:A:2:LYS:N	2:A:410:HOH:O	2.39	0.54
1:A:217:LEU:N	1:A:217:LEU:CD2	2.71	0.53
1:A:90:GLY:HA2	1:A:330:TYR:H	1.73	0.53
1:B:212:ALA:O	1:B:213:GLU:HB2	2.08	0.53
1:C:173:TYR:OH	1:C:184:GLU:OE2	2.19	0.53
1:D:176:ARG:HG2	1:D:176:ARG:NH1	2.22	0.53
1:A:123:ASP:O	1:A:127:ARG:HG2	2.10	0.52
1:A:217:LEU:HD22	1:A:217:LEU:H	1.72	0.52
1:D:237:ARG:NH2	2:D:403:HOH:O	2.24	0.51
1:B:198:LEU:O	1:B:226:MET:HA	2.11	0.51
1:B:328:GLN:HG2	1:D:17:GLU:HB3	1.91	0.51
1:B:193:THR:OG1	1:B:194:LEU:N	2.42	0.51
1:B:142:GLY:O	1:B:146:LYS:HG3	2.11	0.51
1:B:273:ARG:HD2	1:B:274:PRO:O	2.10	0.50
1:A:216:HIS:NE2	1:A:264:GLU:OE2	2.38	0.50
1:C:151:ARG:HD2	1:C:217:LEU:HD22	1.94	0.49
1:A:206:PRO:HD2	1:A:233:ILE:O	2.13	0.49
1:B:173:TYR:HD2	1:B:189:VAL:HG22	1.78	0.49
1:C:153:MET:HG3	1:C:179:LEU:HD11	1.95	0.48
1:D:57:ILE:HG21	1:D:71:ILE:HD13	1.97	0.47
1:A:193:THR:OG1	1:A:196:GLU:HG3	2.14	0.47
1:A:135:ARG:NH1	2:A:421:HOH:O	2.47	0.47
1:C:178:PRO:HG3	1:C:191:ARG:NH1	2.29	0.47
1:A:198:LEU:O	1:A:226:MET:HA	2.15	0.47
1:A:221:THR:O	1:A:225:LYS:HG3	2.15	0.47
1:B:59:SER:HA	1:B:62:LEU:HD12	1.97	0.46
1:D:151:ARG:HB3	1:D:206:PRO:HA	1.97	0.46
1:D:237:ARG:HG2	1:D:240:VAL:HG23	1.97	0.46
1:C:273:ARG:NH2	1:C:275:GLN:HE21	1.98	0.46
1:B:215:LEU:HD22	1:B:270:ARG:HE	1.80	0.46
1:C:194:LEU:HD22	1:C:217:LEU:HD11	1.97	0.46
1:D:322:PRO:HG2	1:D:325:ALA:HB2	1.98	0.46
1:C:57:ILE:HG21	1:C:71:ILE:HD13	1.97	0.46
1:D:14:GLU:OE1	2:D:401:HOH:O	2.21	0.46
1:A:99:LEU:HD12	1:A:305:ILE:HG13	1.97	0.45
1:D:239:SER:CB	1:D:270:ARG:HH22	2.29	0.45
1:B:193:THR:HG22	1:B:196:GLU:OE2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:GLU:OE2	2:A:403:HOH:O	2.21	0.45
1:C:244:ASN:HA	1:C:247:ILE:HD12	1.97	0.45
1:D:258:TYR:O	1:D:288:THR:HA	2.17	0.45
1:D:206:PRO:HD2	1:D:233:ILE:O	2.17	0.45
1:A:100:LEU:HG	1:A:296:SER:HB3	1.98	0.45
1:A:120:LEU:HD11	1:A:289:PHE:CD1	2.52	0.45
1:B:106:GLU:OE2	2:B:402:HOH:O	2.21	0.45
1:C:171:LEU:O	1:C:189:VAL:HA	2.17	0.45
1:A:120:LEU:HD21	1:A:289:PHE:CE1	2.52	0.44
1:C:68:LEU:O	1:C:91:VAL:HG22	2.17	0.44
1:C:198:LEU:O	1:C:226:MET:HA	2.17	0.44
1:C:147:THR:HG22	1:C:201:CYS:HA	2.00	0.44
1:A:139:TYR:CZ	1:D:302:ARG:HG3	2.53	0.43
1:C:236:CYS:SG	1:C:237:ARG:N	2.91	0.43
1:B:100:LEU:HD22	1:B:296:SER:CB	2.48	0.43
1:B:123:ASP:O	1:B:127:ARG:HG2	2.18	0.43
1:B:80:ASN:HB2	2:B:493:HOH:O	2.17	0.43
1:B:258:TYR:O	1:B:288:THR:HA	2.19	0.43
1:D:198:LEU:O	1:D:226:MET:HA	2.19	0.42
1:B:298:VAL:HG23	1:B:301:VAL:HB	2.01	0.42
1:D:237:ARG:HG2	1:D:240:VAL:CG2	2.49	0.42
1:A:263:PHE:O	1:A:266:GLU:HG2	2.18	0.42
1:C:221:THR:HG22	1:C:225:LYS:HE3	2.02	0.42
1:C:215:LEU:HB2	1:C:216:HIS:CD2	2.50	0.42
1:A:135:ARG:HB2	1:A:137:THR:HG22	2.02	0.42
1:D:153:MET:HE3	1:D:179:LEU:HD21	2.01	0.42
1:A:10:TRP:HB2	1:A:30:THR:O	2.20	0.42
1:C:74:ALA:HB1	1:C:305:ILE:HG22	2.01	0.41
1:A:153:MET:CE	1:A:158:ARG:HG2	2.51	0.41
1:C:254:LYS:HA	1:C:254:LYS:HD3	1.95	0.41
1:B:267:GLU:O	1:B:273:ARG:HG3	2.21	0.41
1:D:84:ASN:O	1:D:88:ARG:HG3	2.20	0.41
1:A:146:LYS:HD3	1:A:146:LYS:HA	1.84	0.41
1:D:153:MET:CE	1:D:179:LEU:HD11	2.51	0.41
1:A:47:ASP:OD1	2:A:404:HOH:O	2.22	0.41
1:A:281:LEU:HA	1:A:281:LEU:HD23	1.84	0.41
1:B:180:ASN:OD1	1:B:180:ASN:C	2.59	0.41
1:D:153:MET:HE1	1:D:158:ARG:HA	2.03	0.41
1:C:267:GLU:HB3	1:C:270:ARG:HG2	2.03	0.41
1:D:329:PRO:O	1:D:330:TYR:HB2	2.21	0.41
1:C:151:ARG:CD	1:C:217:LEU:HD22	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:217:LEU:CD2	1:A:217:LEU:H	2.32	0.40
1:B:90:GLY:HA2	1:B:329:PRO:HB2	2.03	0.40
1:D:268:TRP:CD2	1:D:273:ARG:NH1	2.88	0.40
1:C:247:ILE:HD11	1:C:278:PRO:HB2	2.03	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	327/338 (97%)	317 (97%)	10 (3%)	0	100	100
1	B	327/338 (97%)	315 (96%)	11 (3%)	1 (0%)	41	46
1	C	328/338 (97%)	315 (96%)	12 (4%)	1 (0%)	41	46
1	D	322/338 (95%)	315 (98%)	6 (2%)	1 (0%)	41	46
All	All	1304/1352 (96%)	1262 (97%)	39 (3%)	3 (0%)	47	55

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	213	GLU
1	C	180	ASN
1	D	269	ILE

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	264/270 (98%)	263 (100%)	1 (0%)	91	94
1	B	264/270 (98%)	262 (99%)	2 (1%)	81	88
1	C	265/270 (98%)	261 (98%)	4 (2%)	65	75
1	D	261/270 (97%)	256 (98%)	5 (2%)	57	66
All	All	1054/1080 (98%)	1042 (99%)	12 (1%)	73	82

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	216	HIS
1	B	237	ARG
1	B	321	LYS
1	C	1	MET
1	C	182	GLU
1	C	270	ARG
1	C	275	GLN
1	D	2	LYS
1	D	44	LYS
1	D	59	SER
1	D	191	ARG
1	D	217	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	125	GLN
1	C	216	HIS
1	C	275	GLN
1	D	190	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	329/338 (97%)	0.03	14 (4%) 35 37	18, 31, 58, 110	0
1	B	329/338 (97%)	0.18	19 (5%) 23 25	19, 36, 73, 101	0
1	C	330/338 (97%)	0.13	21 (6%) 19 21	19, 35, 71, 112	0
1	D	326/338 (96%)	-0.13	10 (3%) 49 52	19, 31, 57, 109	0
All	All	1314/1352 (97%)	0.05	64 (4%) 29 32	18, 33, 69, 112	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	215	LEU	18.2
1	A	215	LEU	14.8
1	A	269	ILE	13.7
1	A	272	ASP	12.1
1	A	268	TRP	10.7
1	C	271	ALA	10.2
1	C	269	ILE	10.1
1	C	215	LEU	9.9
1	A	271	ALA	9.9
1	C	270	ARG	9.7
1	B	268	TRP	9.4
1	B	272	ASP	9.2
1	B	271	ALA	8.8
1	C	272	ASP	8.6
1	B	269	ILE	8.2
1	B	212	ALA	8.1
1	D	272	ASP	8.1
1	C	216	HIS	7.6
1	C	212	ALA	7.6
1	D	271	ALA	7.4
1	D	269	ILE	7.3

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	C	214	THR	7.0
1	B	211	ALA	6.5
1	A	270	ARG	6.5
1	A	274	PRO	6.2
1	B	330	TYR	5.7
1	C	1	MET	5.7
1	A	273	ARG	5.6
1	D	268	TRP	5.4
1	C	268	TRP	5.3
1	A	213	GLU	5.1
1	A	216	HIS	5.1
1	B	270	ARG	5.0
1	A	330	TYR	4.9
1	C	211	ALA	4.8
1	C	213	GLU	4.5
1	B	214	THR	4.4
1	D	330	TYR	4.4
1	D	270	ARG	4.3
1	C	275	GLN	4.0
1	B	213	GLU	3.8
1	A	275	GLN	3.8
1	B	209	PRO	3.5
1	C	330	TYR	3.4
1	B	273	ARG	3.4
1	A	214	THR	3.3
1	D	216	HIS	3.2
1	B	274	PRO	3.0
1	D	212	ALA	2.9
1	D	211	ALA	2.8
1	C	185	LYS	2.7
1	B	275	GLN	2.7
1	C	217	LEU	2.6
1	D	50	MET	2.6
1	C	188	HIS	2.5
1	B	210	MET	2.4
1	C	181	ALA	2.3
1	C	221	THR	2.3
1	B	260	ALA	2.2
1	B	216	HIS	2.2
1	B	130	HIS	2.2
1	C	273	ARG	2.1
1	A	251	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	176	ARG	2.1

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.